

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102124\
 Data File : VX043457.D
 Acq On : 21 Oct 2024 16:12
 Operator : JC/MD
 Sample : P4419-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :

Quant Time: Oct 22 02:40:45 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 02 16:50:57 2024
 Response via : Initial Calibration

Compound	R.T. QIon		Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.550	168	99072	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	187839	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	167921	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	71876	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	89014	55.015	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	110.020%
35) Dibromofluoromethane	5.391	113	63591	49.089	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.180%
50) Toluene-d8	8.647	98	223329	49.748	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.500%
62) 4-Bromofluorobenzene	11.079	95	83160	50.990	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	101.980%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	37900	56.091	ug/l	97
18) Methyl Acetate	2.709	43	3998	2.359	ug/l #	89
20) Methylene Chloride	2.788	84	3383	2.715	ug/l #	92
43) Isopropyl Acetate	6.348	43	104774	32.277	ug/l	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102124\
Data File : VX043457.D
Acq On : 21 Oct 2024 16:12
Operator : JC/MD
Sample : P4419-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :

Quant Time: Oct 22 02:40:45 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 02 16:50:57 2024
Response via : Initial Calibration

